



Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 01:41 am BST

PDB ID : 5FBW
Title : PI4KB in complex with Rab11 and the MI369 Inhibitor
Authors : Chalupska, D.; Mejdrova, I.; Nencka, R.; Boura, E.
Deposited on : 2015-12-14
Resolution : 3.49 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

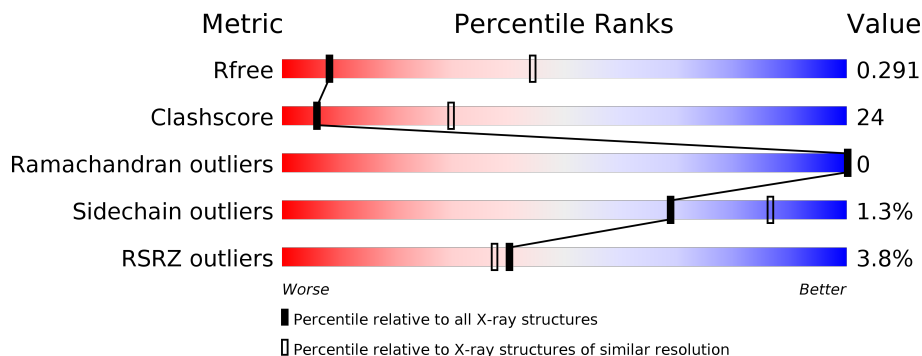
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.49 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1379 (3.56-3.40)
Clashscore	141614	1461 (3.56-3.40)
Ramachandran outliers	138981	1424 (3.56-3.40)
Sidechain outliers	138945	1425 (3.56-3.40)
RSRZ outliers	127900	1289 (3.56-3.40)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	572	 3% 55% 24% 19%
2	B	221	 4% 45% 29% 25%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5037 atoms, of which 14 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Phosphatidylinositol 4-kinase beta, Phosphatidylinositol 4-kinase beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	463	3719	2394	642	659	24	0	0	0

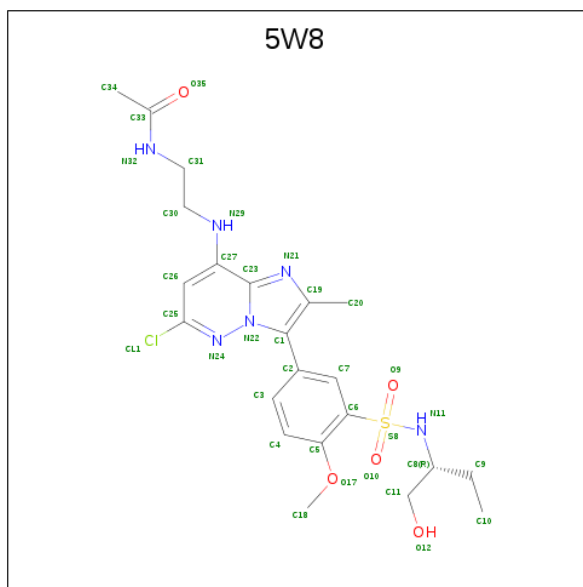
- Molecule 2 is a protein called Ras-related protein Rab-11A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	165	1237	786	207	243	1	0	0	0

There are 6 discrepancies between the modelled and reference sequences:

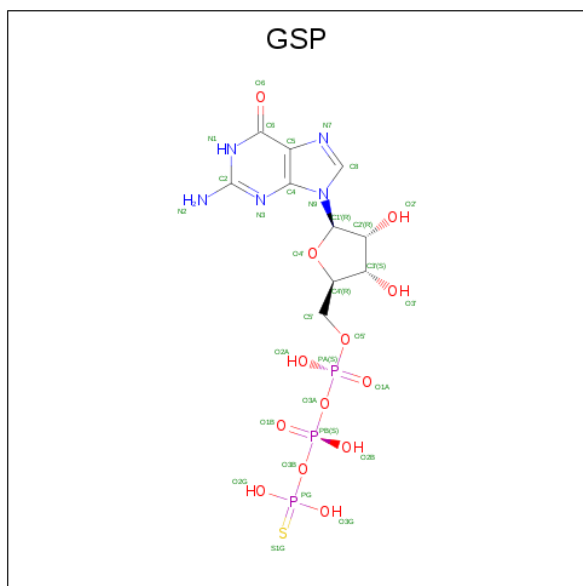
Chain	Residue	Modelled	Actual	Comment	Reference
B	-4	GLY	-	expression tag	UNP P62491
B	-3	ALA	-	expression tag	UNP P62491
B	-2	MET	-	expression tag	UNP P62491
B	-1	GLY	-	expression tag	UNP P62491
B	0	SER	-	expression tag	UNP P62491
B	70	LEU	GLN	engineered mutation	UNP P62491

- Molecule 3 is {N}-[2-[6-chloranyl-3-[4-methoxy-3-[(2 {R})-1-oxidanylbutan-2-yl]sulfamoyl]phenyl]-2-methyl-imidazo[1,2-b]pyridazin-8-yl]amino]ethyl]ethanamide (three-letter code: 5W8) (formula: C₂₂H₂₉ClN₆O₅S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
			Total	C	Cl	N	O			S
3	A	1	35	22	1	6	5	1	0	0

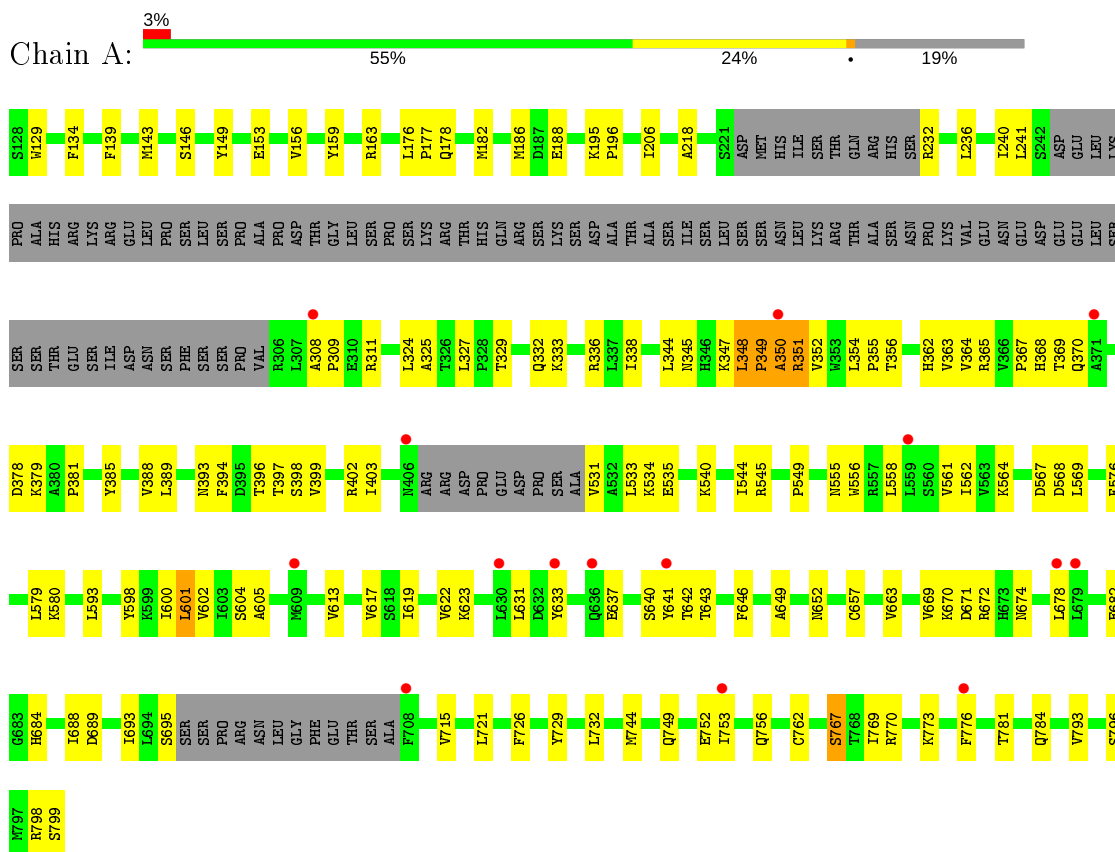
- Molecule 4 is 5'-GUANOSINE-DIPHOSPHATE-MONOTHIOPHOSPHATE (three-letter code: GSP) (formula: C₁₀H₁₆N₅O₁₃P₃S).



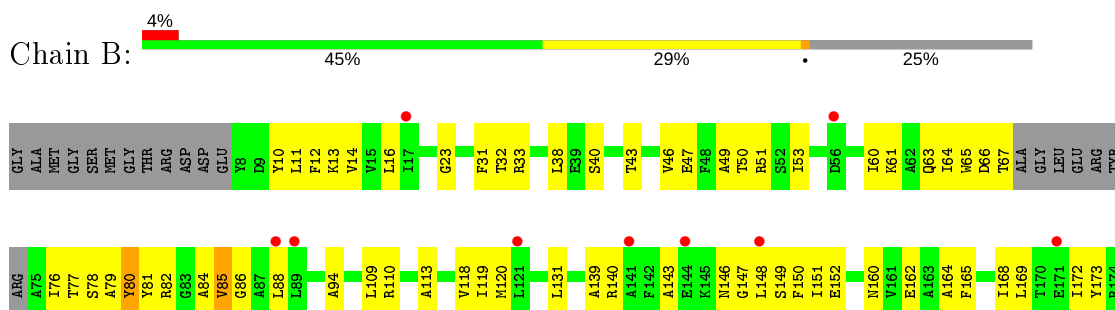
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: Phosphatidylinositol 4-kinase beta, Phosphatidylinositol 4-kinase beta



- Molecule 2: Ras-related protein Rab-11A



I175	V176	Q178	K179	GLN	MET	SER	ASP	ARG	ARG	GLU	ASN	ASP	MET	SER	SER	PRO	SER	ASN	ASN	VAL	VAL	PRO	ILE	ILE	HIS	VAL	VAL	PRO	PRO	THR	THR	THR	GLU	ASN	LYS	PRO	LYS	LYS	VAL	GLN	CYS	CYS	GLN	GLN	ASN	ASN	ILE
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4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	49.07Å 104.09Å 187.56Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.47 – 3.49 47.47 – 3.48	Depositor EDS
% Data completeness (in resolution range)	98.0 (47.47-3.49) 84.8 (47.47-3.48)	Depositor EDS
R_{merge}	0.19	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.62 (at 3.48Å)	Xtrriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.246 , 0.283 0.250 , 0.291	Depositor DCC
R_{free} test set	635 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	104.5	Xtrriage
Anisotropy	0.602	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 85.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.29$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	5037	wwPDB-VP
Average B, all atoms (Å ²)	127.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.11% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GSP, 5W8

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.29	0/3796	0.54	1/5127 (0.0%)
2	B	0.33	0/1256	0.63	0/1712
All	All	0.30	0/5052	0.57	1/6839 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
2	B	0	1
All	All	0	3

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	350	ALA	N-CA-C	5.58	126.08	111.00

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	349	PRO	Peptide
1	A	767	SER	Peptide
2	B	47	GLU	Peptide

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3719	0	3766	161	0
2	B	1237	0	1174	84	0
3	A	35	0	0	8	0
4	B	32	14	12	3	0
All	All	5023	14	4952	244	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 24.

All (244) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:348:LEU:CD1	1:A:365:ARG:HA	1.77	1.14
1:A:348:LEU:HD12	1:A:364:VAL:O	1.49	1.11
2:B:143:ALA:HA	2:B:148:LEU:HD11	1.34	1.05
1:A:348:LEU:HD12	1:A:365:ARG:HA	1.38	1.04
2:B:81:TYR:O	2:B:84:ALA:HB2	1.58	1.03
1:A:564:LYS:HE3	1:A:567:ASP:HB2	1.41	1.03
1:A:345:ASN:OD1	1:A:348:LEU:HG	1.64	0.97
1:A:345:ASN:HA	1:A:348:LEU:HD23	1.45	0.97
1:A:348:LEU:HD12	1:A:364:VAL:C	1.86	0.96
1:A:348:LEU:CD1	1:A:365:ARG:CA	2.45	0.95
2:B:32:THR:HG21	2:B:49:ALA:HB1	1.48	0.93
1:A:562:ILE:HG21	3:A:801:5W8:C10	1.99	0.93
1:A:345:ASN:HA	1:A:348:LEU:CG	1.99	0.92
1:A:139:PHE:HA	1:A:143:MET:HE2	1.51	0.91
1:A:345:ASN:HA	1:A:348:LEU:CD2	2.01	0.90
2:B:81:TYR:O	2:B:84:ALA:CB	2.19	0.89
2:B:140:ARG:HH11	2:B:140:ARG:HB2	1.38	0.88
1:A:345:ASN:HA	1:A:348:LEU:HB2	1.57	0.86
1:A:348:LEU:HD12	1:A:365:ARG:CA	2.03	0.85
1:A:345:ASN:CA	1:A:348:LEU:HD23	2.06	0.83
1:A:348:LEU:HD11	1:A:365:ARG:HA	1.60	0.83
1:A:345:ASN:HA	1:A:348:LEU:CB	2.08	0.83
1:A:348:LEU:HD11	1:A:365:ARG:CA	2.07	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:120:MET:CE	2:B:168:ILE:HB	2.10	0.81
1:A:393:ASN:HB3	1:A:396:THR:HG22	1.63	0.80
1:A:345:ASN:O	1:A:348:LEU:HB2	1.81	0.79
1:A:643:THR:HG23	1:A:646:PHE:H	1.46	0.79
1:A:564:LYS:HD3	1:A:569:LEU:HD21	1.64	0.78
1:A:347:LYS:C	1:A:349:PRO:HD2	2.04	0.77
1:A:345:ASN:CA	1:A:348:LEU:HB2	2.15	0.77
1:A:329:THR:HG23	1:A:332:GLN:H	1.50	0.77
1:A:344:LEU:C	1:A:348:LEU:HD23	2.04	0.76
2:B:140:ARG:HB2	2:B:140:ARG:NH1	2.01	0.76
2:B:12:PHE:HE2	2:B:60:ILE:HD13	1.49	0.76
1:A:601:LEU:HD12	1:A:602:VAL:N	2.02	0.74
1:A:348:LEU:HD12	1:A:365:ARG:N	2.02	0.74
1:A:345:ASN:N	1:A:348:LEU:HD23	2.02	0.73
1:A:393:ASN:HB3	1:A:396:THR:CG2	2.17	0.73
1:A:348:LEU:CD1	1:A:364:VAL:C	2.57	0.73
1:A:348:LEU:CD1	1:A:365:ARG:N	2.52	0.73
2:B:140:ARG:HA	2:B:143:ALA:HB3	1.69	0.73
2:B:23:GLY:N	4:B:301:GSP:O2B	2.21	0.72
1:A:370:GLN:NE2	1:A:531:VAL:O	2.18	0.71
1:A:579:LEU:HD21	1:A:663:VAL:HG11	1.72	0.71
1:A:403:ILE:HG22	1:A:549:PRO:HB3	1.72	0.71
1:A:564:LYS:CE	1:A:567:ASP:HB2	2.20	0.71
2:B:60:ILE:HG22	2:B:61:LYS:N	2.06	0.71
1:A:564:LYS:HE2	1:A:569:LEU:HD21	1.71	0.70
1:A:715:VAL:HG11	1:A:721:LEU:HD23	1.73	0.70
1:A:402:ARG:O	1:A:403:ILE:HD13	1.90	0.69
1:A:348:LEU:N	1:A:349:PRO:CD	2.54	0.69
1:A:593:LEU:HD21	1:A:732:LEU:HB3	1.75	0.69
1:A:327:LEU:HD11	1:A:336:ARG:NE	2.07	0.69
2:B:143:ALA:CA	2:B:148:LEU:HD11	2.17	0.68
1:A:753:ILE:O	1:A:756:GLN:HB2	1.94	0.68
2:B:175:ILE:HD12	2:B:175:ILE:C	2.13	0.68
2:B:140:ARG:HA	2:B:143:ALA:CB	2.23	0.67
1:A:348:LEU:CD1	1:A:364:VAL:O	2.35	0.67
2:B:120:MET:HA	2:B:149:SER:HB2	1.75	0.67
2:B:172:ILE:O	2:B:175:ILE:HG23	1.95	0.67
1:A:351:ARG:HA	1:A:362:HIS:HD2	1.59	0.67
2:B:65:TRP:CZ3	2:B:80:TYR:CE1	2.83	0.66
2:B:32:THR:HG21	2:B:49:ALA:CB	2.25	0.66
2:B:94:ALA:HB1	2:B:131:LEU:CD2	2.26	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:350:ALA:HB1	1:A:352:VAL:HG22	1.77	0.65
2:B:120:MET:HE3	2:B:168:ILE:HB	1.78	0.65
1:A:564:LYS:CD	1:A:569:LEU:HD21	2.25	0.64
2:B:13:LYS:HG3	2:B:84:ALA:HA	1.78	0.64
1:A:345:ASN:C	1:A:348:LEU:HB2	2.16	0.64
1:A:218:ALA:HB1	1:A:605:ALA:HB2	1.80	0.64
2:B:16:LEU:O	2:B:67:THR:HG22	1.97	0.64
2:B:12:PHE:CE2	2:B:60:ILE:HD13	2.33	0.64
2:B:120:MET:CB	2:B:149:SER:HB2	2.27	0.63
1:A:327:LEU:HD11	1:A:336:ARG:CZ	2.28	0.63
1:A:652:ASN:HB3	1:A:684:HIS:CD2	2.34	0.63
1:A:329:THR:HG22	1:A:332:GLN:HB2	1.81	0.62
1:A:669:VAL:HG22	1:A:693:ILE:HG22	1.79	0.62
1:A:347:LYS:C	1:A:349:PRO:CD	2.67	0.62
1:A:206:ILE:HD11	1:A:394:PHE:CD2	2.35	0.61
1:A:715:VAL:HG11	1:A:721:LEU:CD2	2.30	0.61
1:A:593:LEU:CD2	1:A:732:LEU:HB3	2.30	0.61
2:B:65:TRP:CZ3	2:B:80:TYR:HE1	2.19	0.61
2:B:143:ALA:HA	2:B:148:LEU:CD1	2.21	0.61
2:B:66:ASP:OD1	2:B:67:THR:N	2.33	0.61
1:A:348:LEU:N	1:A:349:PRO:HD3	2.16	0.60
2:B:109:LEU:O	2:B:113:ALA:HB3	2.01	0.60
1:A:134:PHE:O	1:A:163:ARG:NH1	2.34	0.60
1:A:351:ARG:HB2	1:A:399:VAL:HG11	1.81	0.60
1:A:767:SER:O	1:A:770:ARG:N	2.34	0.60
1:A:355:PRO:HG2	1:A:601:LEU:CD2	2.32	0.60
1:A:188:GLU:HG3	1:A:232:ARG:HH12	1.67	0.60
1:A:241:LEU:HD22	1:A:311:ARG:HD2	1.82	0.60
1:A:672:ARG:NH2	1:A:689:ASP:O	2.35	0.60
2:B:131:LEU:O	2:B:131:LEU:HD23	2.02	0.59
2:B:12:PHE:CE2	2:B:60:ILE:HG21	2.38	0.59
1:A:564:LYS:CE	1:A:569:LEU:HD21	2.33	0.58
2:B:60:ILE:HG22	2:B:61:LYS:H	1.67	0.58
1:A:598:TYR:CE1	1:A:600:ILE:HD11	2.39	0.58
1:A:564:LYS:HG2	3:A:801:5W8:O12	2.04	0.58
2:B:77:THR:HG23	2:B:77:THR:O	2.02	0.58
1:A:347:LYS:O	1:A:349:PRO:HD2	2.02	0.58
2:B:120:MET:HB2	2:B:149:SER:HB2	1.84	0.58
2:B:109:LEU:CD1	2:B:119:ILE:HD11	2.34	0.57
2:B:51:ARG:NH2	2:B:162:GLU:OE2	2.33	0.57
1:A:356:THR:HG21	1:A:561:VAL:CG1	2.34	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:120:MET:HE1	2:B:168:ILE:HB	1.85	0.57
2:B:50:THR:HG22	2:B:63:GLN:HA	1.86	0.57
1:A:195:LYS:HB3	1:A:196:PRO:HD3	1.86	0.57
1:A:564:LYS:HZ2	3:A:801:5W8:C11	2.18	0.57
2:B:151:ILE:HD11	2:B:160:ASN:HB3	1.86	0.57
1:A:744:MET:CE	1:A:773:LYS:HA	2.34	0.56
1:A:355:PRO:HG2	1:A:601:LEU:HD22	1.87	0.56
1:A:381:PRO:HG3	3:A:801:5W8:C11	2.36	0.56
1:A:641:TYR:CE1	1:A:642:THR:HG23	2.41	0.56
1:A:576:PHE:CD1	1:A:598:TYR:O	2.59	0.56
1:A:640:SER:O	1:A:643:THR:HG22	2.06	0.56
1:A:613:VAL:O	3:A:801:5W8:N29	2.39	0.55
1:A:363:VAL:HA	1:A:388:VAL:HG12	1.88	0.55
1:A:564:LYS:NZ	3:A:801:5W8:C11	2.69	0.55
2:B:13:LYS:CG	2:B:84:ALA:HA	2.37	0.55
1:A:752:GLU:HA	1:A:769:ILE:HD11	1.89	0.55
1:A:356:THR:HG21	1:A:561:VAL:HG12	1.89	0.54
2:B:12:PHE:CD2	2:B:60:ILE:HG21	2.43	0.54
1:A:657:CYS:HB3	1:A:729:TYR:HE1	1.72	0.54
2:B:120:MET:CA	2:B:149:SER:HB2	2.37	0.54
1:A:351:ARG:HA	1:A:362:HIS:CD2	2.42	0.54
2:B:152:GLU:O	2:B:160:ASN:ND2	2.32	0.54
1:A:695:SER:HB2	1:A:762:CYS:O	2.08	0.54
2:B:165:PHE:O	2:B:168:ILE:HG22	2.08	0.54
1:A:397:THR:HG22	1:A:398:SER:N	2.22	0.53
1:A:545:ARG:HG3	1:A:556:TRP:CZ3	2.43	0.53
1:A:641:TYR:CD1	1:A:641:TYR:C	2.81	0.53
2:B:60:ILE:CG2	2:B:61:LYS:N	2.70	0.53
2:B:175:ILE:O	2:B:175:ILE:HD12	2.09	0.53
1:A:344:LEU:O	1:A:348:LEU:HD23	2.08	0.53
1:A:403:ILE:CG2	1:A:549:PRO:HB3	2.39	0.53
2:B:14:VAL:HA	2:B:86:GLY:O	2.09	0.52
1:A:767:SER:O	1:A:770:ARG:HB2	2.09	0.52
2:B:140:ARG:CA	2:B:143:ALA:HB3	2.39	0.52
1:A:149:TYR:HB2	1:A:182:MET:CE	2.40	0.52
1:A:535:GLU:HG2	1:A:540:LYS:HE3	1.90	0.52
1:A:350:ALA:O	1:A:351:ARG:HB3	2.10	0.52
2:B:120:MET:HE1	2:B:164:ALA:O	2.08	0.52
1:A:308:ALA:HB3	1:A:309:PRO:HD3	1.91	0.52
1:A:657:CYS:HB3	1:A:729:TYR:CE1	2.45	0.52
2:B:140:ARG:O	2:B:143:ALA:HB3	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:329:THR:HG22	1:A:332:GLN:CG	2.40	0.51
1:A:619:ILE:HB	1:A:674:ASN:HB3	1.91	0.51
2:B:81:TYR:O	2:B:82:ARG:C	2.46	0.51
2:B:109:LEU:HD11	2:B:119:ILE:HD11	1.92	0.51
1:A:402:ARG:C	1:A:403:ILE:HD13	2.31	0.51
1:A:568:ASP:C	1:A:569:LEU:HD23	2.30	0.51
1:A:378:ASP:OD1	1:A:379:LYS:N	2.44	0.51
1:A:533:LEU:HG	1:A:534:LYS:N	2.26	0.51
1:A:617:VAL:HG23	1:A:622:VAL:HG23	1.93	0.51
2:B:94:ALA:HB1	2:B:131:LEU:HD23	1.93	0.51
2:B:60:ILE:CG2	2:B:61:LYS:H	2.24	0.50
1:A:348:LEU:HD11	1:A:365:ARG:C	2.32	0.50
2:B:76:ILE:HG13	2:B:77:THR:N	2.26	0.50
1:A:544:ILE:HD12	1:A:558:LEU:CD2	2.42	0.50
1:A:744:MET:HE3	1:A:776:PHE:CD1	2.47	0.50
1:A:781:THR:HG23	1:A:784:GLN:H	1.77	0.50
1:A:329:THR:HG22	1:A:332:GLN:CB	2.42	0.50
1:A:325:ALA:HA	1:A:378:ASP:O	2.12	0.50
1:A:368:HIS:CE1	1:A:369:THR:HG23	2.48	0.49
2:B:151:ILE:CD1	2:B:160:ASN:HB3	2.42	0.49
1:A:726:PHE:O	1:A:729:TYR:HB3	2.13	0.49
2:B:65:TRP:HZ3	2:B:80:TYR:CE1	2.28	0.49
2:B:146:ASN:OD1	2:B:147:GLY:N	2.42	0.49
1:A:236:LEU:O	1:A:240:ILE:HG13	2.13	0.48
1:A:533:LEU:HG	1:A:534:LYS:H	1.78	0.48
2:B:10:TYR:O	2:B:11:LEU:HD12	2.12	0.48
2:B:10:TYR:C	2:B:11:LEU:HD12	2.34	0.48
1:A:153:GLU:OE1	2:B:40:SER:OG	2.22	0.48
1:A:544:ILE:HD12	1:A:558:LEU:HD21	1.96	0.48
1:A:324:LEU:O	1:A:333:LYS:HE3	2.14	0.48
1:A:535:GLU:OE2	1:A:540:LYS:HE2	2.13	0.48
1:A:604:SER:O	1:A:605:ALA:HB3	2.13	0.48
2:B:109:LEU:C	2:B:109:LEU:HD23	2.35	0.47
1:A:308:ALA:N	1:A:309:PRO:CD	2.77	0.47
1:A:678:LEU:HD22	1:A:688:ILE:HG21	1.95	0.47
1:A:670:LYS:HA	1:A:670:LYS:HD3	1.69	0.47
1:A:139:PHE:HD1	1:A:143:MET:CE	2.28	0.47
1:A:129:TRP:N	1:A:129:TRP:CD1	2.83	0.46
1:A:389:LEU:HB3	1:A:555:ASN:OD1	2.15	0.46
2:B:53:ILE:HD13	2:B:169:LEU:HD12	1.98	0.46
1:A:752:GLU:HA	1:A:769:ILE:CD1	2.45	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:156:VAL:HG22	2:B:38:LEU:HB3	1.98	0.45
1:A:329:THR:CG2	1:A:332:GLN:H	2.25	0.45
1:A:176:LEU:N	1:A:177:PRO:CD	2.79	0.45
2:B:109:LEU:HD13	2:B:119:ILE:HD11	1.98	0.45
1:A:564:LYS:HE3	1:A:567:ASP:CB	2.28	0.45
1:A:793:VAL:O	1:A:796:SER:OG	2.24	0.45
2:B:43:THR:HB	4:B:301:GSP:S1G	2.56	0.45
1:A:649:ALA:HA	1:A:682:GLU:O	2.16	0.45
1:A:345:ASN:HA	1:A:348:LEU:HG	1.91	0.45
1:A:598:TYR:OH	1:A:688:ILE:O	2.25	0.45
1:A:619:ILE:O	1:A:623:LYS:HG3	2.16	0.44
1:A:329:THR:CG2	1:A:332:GLN:HG3	2.47	0.44
1:A:381:PRO:HB3	3:A:801:5W8:O12	2.17	0.44
1:A:350:ALA:HB1	1:A:352:VAL:CG2	2.45	0.44
2:B:118:VAL:HG21	2:B:172:ILE:HD13	2.00	0.44
2:B:168:ILE:O	2:B:172:ILE:HG12	2.17	0.44
2:B:175:ILE:CD1	2:B:175:ILE:C	2.81	0.44
1:A:781:THR:HG22	1:A:784:GLN:CG	2.48	0.44
2:B:43:THR:CB	4:B:301:GSP:S1G	3.06	0.44
3:A:801:5W8:N24	3:A:801:5W8:C7	2.81	0.44
2:B:16:LEU:CD2	2:B:88:LEU:HB2	2.48	0.44
1:A:631:LEU:O	1:A:631:LEU:HD12	2.17	0.43
2:B:173:TYR:O	2:B:175:ILE:N	2.44	0.43
2:B:151:ILE:HG12	2:B:152:GLU:N	2.33	0.43
1:A:393:ASN:CB	1:A:396:THR:HG22	2.41	0.43
1:A:564:LYS:HE2	1:A:569:LEU:CD2	2.45	0.43
2:B:143:ALA:HB1	2:B:150:PHE:HB3	2.01	0.43
1:A:601:LEU:HD12	1:A:602:VAL:H	1.82	0.43
2:B:176:VAL:HG23	2:B:177:SER:N	2.33	0.43
1:A:149:TYR:HB2	1:A:182:MET:HE3	2.01	0.42
2:B:140:ARG:HH11	2:B:140:ARG:CB	2.19	0.42
2:B:85:VAL:HG23	2:B:172:ILE:HD12	2.01	0.42
2:B:109:LEU:HD23	2:B:110:ARG:N	2.34	0.42
2:B:31:PHE:CZ	2:B:165:PHE:CB	3.03	0.42
1:A:159:TYR:OH	1:A:163:ARG:HD2	2.19	0.42
1:A:744:MET:HG3	1:A:776:PHE:CG	2.55	0.42
1:A:564:LYS:CE	1:A:567:ASP:CB	2.93	0.42
1:A:732:LEU:HD23	1:A:732:LEU:HA	1.90	0.42
2:B:53:ILE:CD1	2:B:169:LEU:HD12	2.50	0.42
2:B:139:ALA:O	2:B:143:ALA:N	2.53	0.41
1:A:195:LYS:N	1:A:196:PRO:CD	2.83	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:798:ARG:O	1:A:799:SER:CB	2.68	0.41
2:B:78:SER:OG	2:B:79:ALA:N	2.52	0.41
1:A:146:SER:HB2	1:A:749:GLN:NE2	2.35	0.41
1:A:781:THR:HG22	1:A:784:GLN:HG3	2.02	0.41
1:A:670:LYS:O	1:A:671:ASP:HB3	2.21	0.41
2:B:14:VAL:HB	2:B:64:ILE:HD12	2.02	0.41
1:A:354:LEU:HA	1:A:355:PRO:HD2	1.88	0.41
1:A:633:TYR:CZ	1:A:637:GLU:HG3	2.56	0.41
2:B:32:THR:HG22	2:B:33:ARG:HG3	2.03	0.41
1:A:182:MET:HA	1:A:186:MET:HG3	2.03	0.41
1:A:367:PRO:HD2	1:A:385:TYR:O	2.21	0.41
1:A:672:ARG:NE	1:A:672:ARG:HA	2.36	0.40
2:B:109:LEU:CD1	2:B:119:ILE:CD1	3.00	0.40
1:A:178:GLN:N	1:A:178:GLN:OE1	2.54	0.40
1:A:601:LEU:C	1:A:601:LEU:HD12	2.42	0.40
1:A:338:ILE:HG12	1:A:368:HIS:O	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	453/572 (79%)	439 (97%)	14 (3%)	0	100	100
2	B	161/221 (73%)	152 (94%)	9 (6%)	0	100	100
All	All	614/793 (77%)	591 (96%)	23 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	410/517 (79%)	406 (99%)	4 (1%)	76	89
2	B	125/191 (65%)	122 (98%)	3 (2%)	49	75
All	All	535/708 (76%)	528 (99%)	7 (1%)	69	86

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	348	LEU
1	A	351	ARG
1	A	580	LYS
1	A	601	LEU
2	B	46	VAL
2	B	80	TYR
2	B	85	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry

2 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	GSP	B	301	-	26,34,34	4.22	13 (50%)	28,54,54	2.61	7 (25%)
3	5W8	A	801	-	33,37,37	2.56	6 (18%)	38,53,53	2.68	14 (36%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	GSP	B	301	-	-	8/17/38/38	0/3/3/3
3	5W8	A	801	-	-	17/28/28/28	0/3/3/3

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	801	5W8	C6-S8	-10.61	1.62	1.77
4	B	301	GSP	C2'-C3'	-10.59	1.24	1.53
4	B	301	GSP	O4'-C1'	8.15	1.52	1.41
4	B	301	GSP	O4'-C4'	-7.19	1.28	1.45
3	A	801	5W8	C1-C19	6.99	1.48	1.37
4	B	301	GSP	C6-C5	6.93	1.53	1.41
4	B	301	GSP	C4-N3	6.47	1.45	1.35
4	B	301	GSP	C3'-C4'	5.41	1.66	1.53
4	B	301	GSP	C2-N2	5.05	1.44	1.33
4	B	301	GSP	C6-N1	4.70	1.41	1.33
4	B	301	GSP	C2-N1	4.28	1.43	1.35
3	A	801	5W8	C27-C23	3.90	1.49	1.42
3	A	801	5W8	C23-N21	3.76	1.36	1.33
4	B	301	GSP	O2'-C2'	3.65	1.51	1.43
3	A	801	5W8	C25-N24	3.12	1.36	1.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	301	GSP	C2'-C1'	3.07	1.58	1.53
3	A	801	5W8	C25-CL1	2.95	1.80	1.74
4	B	301	GSP	C2-N3	2.76	1.47	1.34
4	B	301	GSP	PG-S1G	2.41	1.95	1.90

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	301	GSP	C1'-N9-C4	10.09	144.37	126.64
3	A	801	5W8	O10-S8-O9	-8.58	109.00	119.55
3	A	801	5W8	O17-C5-C6	6.12	121.25	116.50
3	A	801	5W8	C8-N11-S8	-5.80	112.04	122.08
4	B	301	GSP	N3-C2-N1	-5.36	120.07	127.22
3	A	801	5W8	CL1-C25-N24	5.09	120.41	115.19
4	B	301	GSP	C2-N3-C4	3.78	119.67	115.36
3	A	801	5W8	C7-C2-C1	3.69	123.92	119.65
3	A	801	5W8	O17-C5-C4	-3.33	118.67	124.37
3	A	801	5W8	C30-N29-C27	-3.26	117.87	123.28
3	A	801	5W8	C26-C25-N24	-3.19	120.35	125.03
3	A	801	5W8	C19-N21-C23	2.99	108.87	103.59
4	B	301	GSP	C5-C6-N1	-2.61	119.86	123.43
4	B	301	GSP	C6-N1-C2	2.40	119.75	115.93
3	A	801	5W8	O9-S8-C6	2.36	111.55	107.66
3	A	801	5W8	C7-C6-C5	-2.29	117.83	120.04
3	A	801	5W8	O10-S8-C6	2.28	111.41	107.66
4	B	301	GSP	N2-C2-N1	2.22	120.71	117.25
4	B	301	GSP	PA-O3A-PB	-2.22	125.22	132.83
3	A	801	5W8	C25-C26-C27	2.18	120.09	116.85
3	A	801	5W8	C18-O17-C5	2.12	120.74	117.53

There are no chirality outliers.

All (25) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	301	GSP	C5'-O5'-PA-O3A
3	A	801	5W8	C19-C1-C2-C3
3	A	801	5W8	C19-C1-C2-C7
3	A	801	5W8	N22-C1-C2-C3
3	A	801	5W8	N22-C1-C2-C7
3	A	801	5W8	C6-C5-O17-C18
3	A	801	5W8	C5-C6-S8-O9
3	A	801	5W8	C5-C6-S8-O10

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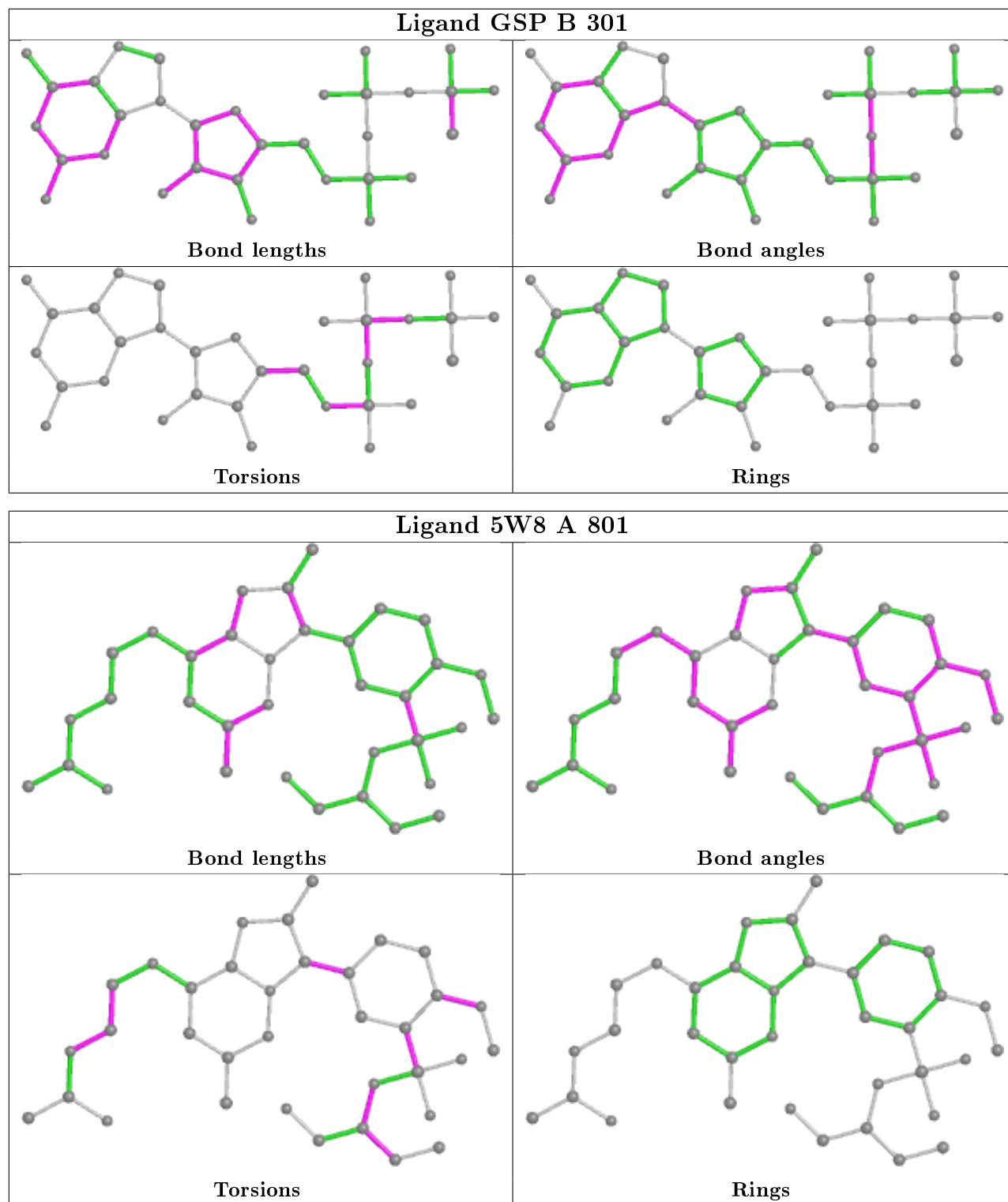
Mol	Chain	Res	Type	Atoms
3	A	801	5W8	C5-C6-S8-N11
3	A	801	5W8	C11-C8-C9-C10
3	A	801	5W8	N11-C8-C9-C10
3	A	801	5W8	C9-C8-N11-S8
3	A	801	5W8	C11-C8-N11-S8
3	A	801	5W8	C30-C31-N32-C33
3	A	801	5W8	N29-C30-C31-N32
3	A	801	5W8	C4-C5-O17-C18
3	A	801	5W8	C7-C6-S8-O9
3	A	801	5W8	C7-C6-S8-N11
4	B	301	GSP	PA-O3A-PB-O3B
4	B	301	GSP	C5'-O5'-PA-O1A
4	B	301	GSP	C5'-O5'-PA-O2A
4	B	301	GSP	O4'-C4'-C5'-O5'
4	B	301	GSP	PG-O3B-PB-O2B
4	B	301	GSP	PA-O3A-PB-O1B
4	B	301	GSP	PA-O3A-PB-O2B

There are no ring outliers.

2 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	301	GSP	3	0
3	A	801	5W8	8	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	463/572 (80%)	0.35	15 (3%) 47 45	91, 124, 147, 161	0
2	B	165/221 (74%)	0.50	9 (5%) 25 24	107, 138, 165, 184	0
All	All	628/793 (79%)	0.39	24 (3%) 40 37	91, 127, 157, 184	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	678	LEU	3.5
1	A	406	ASN	3.5
2	B	88	LEU	3.4
2	B	148	LEU	3.4
1	A	630	LEU	3.3
1	A	708	PHE	3.2
1	A	641	TYR	2.6
1	A	636	GLN	2.6
2	B	89	LEU	2.5
1	A	559	LEU	2.5
2	B	171	GLU	2.3
2	B	17	ILE	2.3
1	A	753	ILE	2.2
1	A	308	ALA	2.2
1	A	776	PHE	2.2
1	A	633	TYR	2.1
2	B	121	LEU	2.1
1	A	679	LEU	2.1
1	A	350	ALA	2.0
1	A	371	ALA	2.0
2	B	144	GLU	2.0
1	A	609	MET	2.0
2	B	56	ASP	2.0
2	B	141	ALA	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

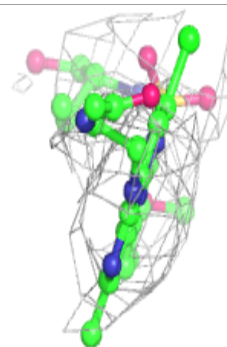
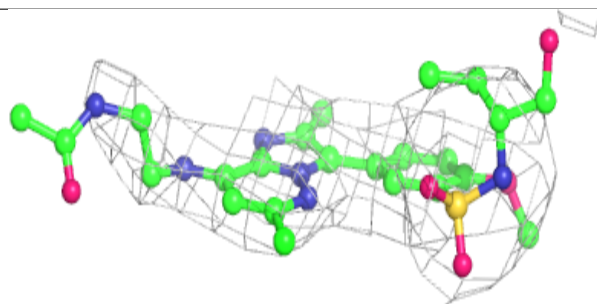
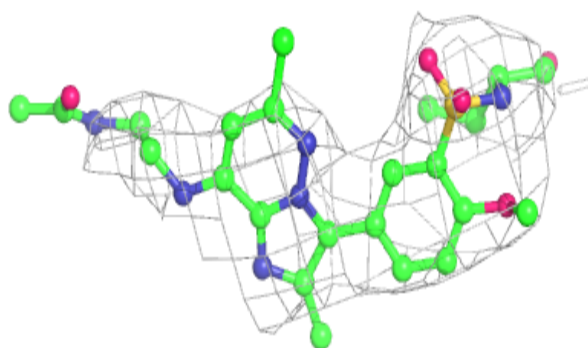
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	5W8	A	801	35/35	0.86	0.35	110,124,139,189	0
4	GSP	B	301	32/32	0.94	0.18	105,120,144,148	0

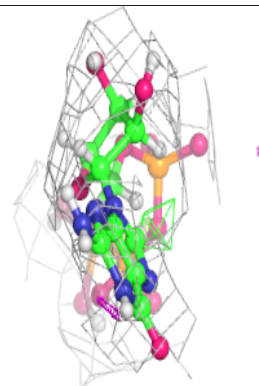
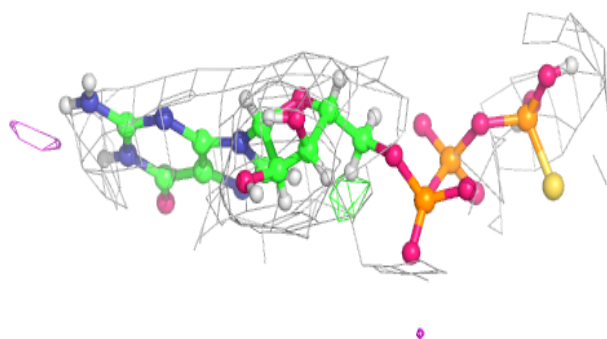
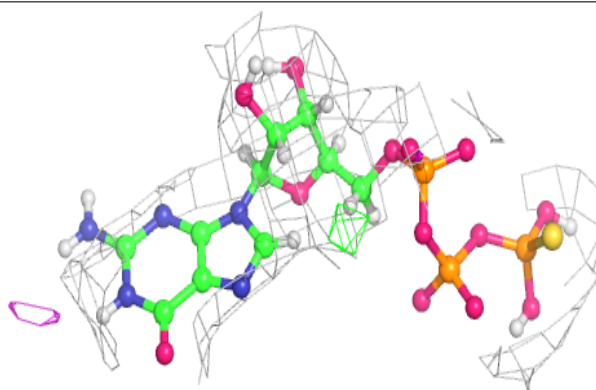
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around 5W8 A 801:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around GSP B 301:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers

There are no such residues in this entry.